Monograph number: 6962.

Title: Oleandomycin.

CAS Registry number: [3922-90-5]

Drug code(s): PA-105;

Trade name(s): Amimycin; Landomycin; Romicil.

Molecular formula: $C_{35}H_{61}NO_{12}$;

Molecular weight: 687.87.

Percent Composition: C 61.11%, H 8.94%, N 2.04%, O 27.91%.

Literature references: Antibiotic substance produced by Streptomyces antibioticus no. ATCC 11891: Sobin et al.; Ratajak, Nubel, U.S. pats. 2,757,123; 2,842,481 (1956, 1958 to Pfizer). Structure: Hochstein et al., J. Am. Chem. Soc. 82, 3225 (1960). Absolute configuration: Celmer, ibid. 87, 1797 (1965); Celmer, Hobbs, Carbohyd. Res. 1, 137 (1965); S. Omura et al., Tetrahedron Letters 1975, 2939. Synthetic study: K. Tatsuta et al., ibid. 29, 3975 (1988). Activity: Hahn, Antibiotics 1, 378, 755 (1967). For a review of macrolide antibiotics see Keller-Schierlein, Fortschr. Chem. Org. Naturst. 30, 313-460 (1973). Toxicity: H. Sous et al., Arzneimittel-Forsch. 8, 386 (1958).

Properties: White amorphous powder. uv max (methanol): 286-289 nm. Moderately sol in water. Sol in dil acids. Freely sol in methanol, ethanol, butanol, acetone. Practically insol in hexane, carbon tetrachloride, dibutyl ether.

Derivative: Hydrochloride,

Molecular Formula: C₃₅H₆₁NO₁₂.HCl,

CAS Registry: [6696-47-5]

Properties: long needles from ethyl acetate, mp 134-135° . $[\alpha]_D^{25}$ –54° (methanol) . Freely sol in water. Forms various cryst hydrates. LD₅₀ in mice, rats (mg/kg): 8200 , >10000 orally ; 600 ,

400 i.v. (Sous).

Melting Point: mp 134-135°

Rotation: -54°

Derivative: Phosphate,

Molecular Formula: C₃₅H₆₁NO₁₂.H₃PO₄,

CAS Registry: [7060-74-4]

Trade name(s): Matromycin (Pfizer).

Derivative: Triacetyl deriv see Troleandomycin.

THERAP CAT: Antibacterial.

THERAP CAT (VET): Antibacterial.

UV Maxima: 286-289